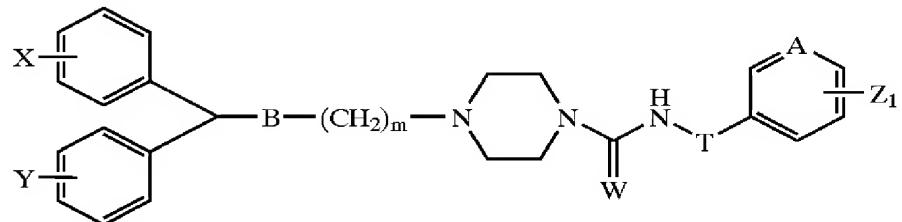


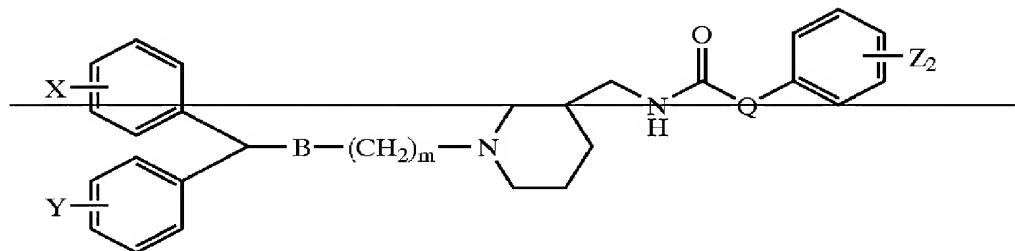
IN THE CLAIMS

1. A compound of Formula I having high affinity for a dopamine transporter having a formula selected from the group consisting of:

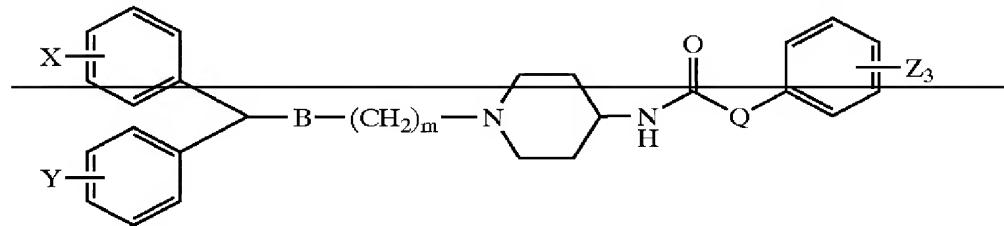
Formula I



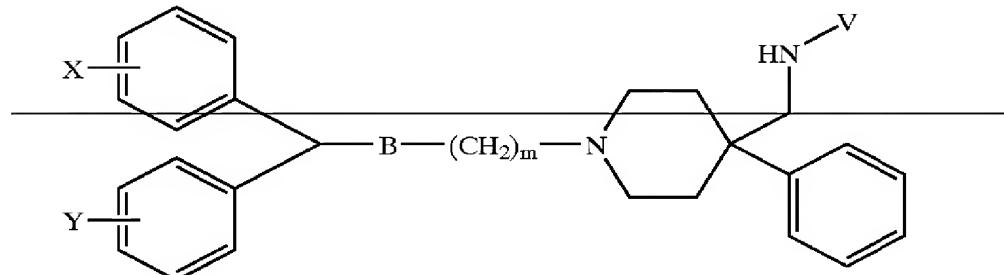
Formula II



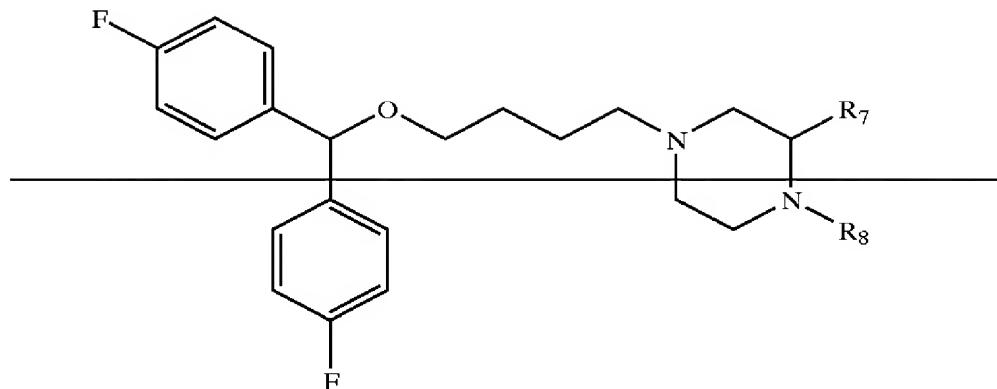
Formula III



Formula IV



Formula V



wherein:

B is -O-, -N(H)-, -C(=O)N(H)-, or -N(H)C(=O)-;

[[n]] m is an integer of 1 to 6; X, Y, Z₄ and Z₂ can be the same or different and are hydrogen, halo, haloalkyl, alkyl, aryl, (C₁-C₆) alkoxy, N-alkyl, (C₂-C₆) acyloxy, N-alkylene, SH, SR, wherein R is from the same group as R₁ and R₂ and can be the same or different than R₁ and R₂, amino, nitro, cyano, hydroxy, C(=O)OR₆, C(=O)NR₅R₆, NR₃R₂ or S(=O)_kR₄ wherein k is 1 or 2, and R₁ to R₆ are independently hydrogen or (C₁-C₆) alkyl;

X is hydrogen, halo, haloalkyl, alkyl, aryl, C₁-C₆ alkoxy, N-alkyl, C₂-C₆ acyloxy, or N-alkylene;

Y is hydrogen, halo, haloalkyl, alkyl, aryl, C₁-C₆ alkoxy, N-alkyl, C₂-C₆ acyloxy, or N-alkylene;

Z₁ is hydrogen, halo, haloalkyl, alkyl, aryl, C₁-C₆ alkoxy, N-alkyl, C₂-C₆ acyloxy, or N-alkylene;

R₁ and R₂ can be the same or different and are hydrogen, (C₁-C₆) alkyl, hydroxyalkyl or mercaptoalkyl, C(=O)OR₄, cyano, (C₁-C₆) alkenyl, (C₂-C₆) alkynyl, or 1,2,4-oxadiazol-5-yl optionally substituted at the 3-position by Z, wherein any (C₁-C₆) alkyl, (C₁-C₆) alkanoyl, (C₂-C₆) alkenyl or (C₂-C₆) alkynyl can optionally be substituted by 1, 2 or 3 Z;

Z₄ is (C₁-C₆) alkyl or phenyl, optionally substituted by 1, 2 or 3 Z₄

~~R₇ can be hydrogen, O or phenyl~~

~~R₈ can be hydrogen, phenyl, halophenyl, nitrophenyl, pyridyl, piperonyl or sulfoxonitrophenyl~~

W is O or S;

T is amino or C₁-C₆ aminoalkyl

A is N or C; and

T is C₁-C₆ alkyl or sulfonyl, and

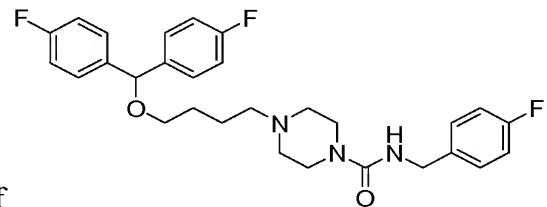
V is alkyl (C₀-C₆), alkenyl, alkynyl, haloaryl, alkyl phenol, alkyl halophenyl, and R₁ or R₂ as indicated above and

Φ is phenyl, naphthyl, thienyl or pyridinyl.

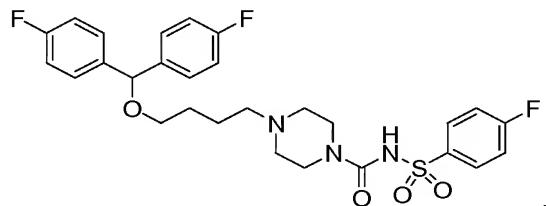
Claims 2-13 (canceled)

14. (new) The compound of claim 1, wherein X is halo.
15. (new) The compound of claim 1, wherein X is *para*-F.
16. (new) The compound of claim 1, wherein Y is halo.
17. (new) The compound of claim 1, wherein Y is *para*-F.
18. (new) The compound of claim 1, wherein B is O.
19. (new) The compound of claim 1, wherein m is 4.
20. (new) The compound of claim 1, wherein W is O.
21. (new) The compound of claim 1, wherein T is C₁-C₆ alkyl.
22. (new) The compound of claim 1, wherein T is -CH₂-.
23. (new) The compound of claim 1, wherein T is sulfonyl.
24. (new) The compound of claim 1, wherein A is C.
25. (new) The compound of claim 1, wherein Z₁ is halo.
26. (new) The compound of claim 1, wherein Z₁ is *para*-F.

27. (new) The compound of claim 1, wherein the compound is selected from the group



consisting of



and

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